Sparse Random Graphs-I

Souvik Dhara

Graph Limits and Processes on Networks: From Epidemics to Misinformation Boot Camp Lets start with a few questions...



How **DISEASE** becomes **Epidemic**?

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What causes INTERNET to breakdown?

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When does **MISINFORMATION** reach a large population?

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➤ Random Graphs provide a *simplified probabilistic representation* to model these complex system.

- ➡ Capture structural properties (degree distribution, communities)
- Provide insight into emergence of different types of behavior such as phase transition

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Random Graphs serve to get provable guarantees for graph algorithms *Example:* Heuristic algorithms for NP-hard problems such as graph partitioning, coloring.

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Plan

Today:

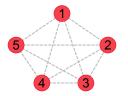
- > Local Branching Process approximation technique on random graphs
- > Explore its relation to *Giant Component Problem* on different models

Tomorrow:

- > Applications to Percolation, Epidemics
- > Using Stochastic Process convergence in Random Graphs

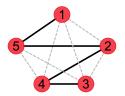
Let's start with the most elementary yet fundamental model...

Erdős-Rényi Random Graph



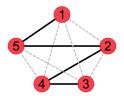
Definition

> Given n nodes $\{1, 2, \ldots, n\}$



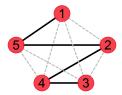
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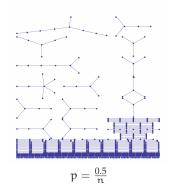
Historical note:

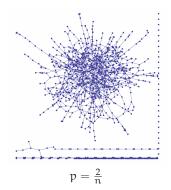
➤ This model was actually studied by Gilbert (1959) and heuristically by Solomonoff & Rapoport (1951)

➤ Erdős & Rényi (1959) initially worked with a slightly different model where fixed number of edges sampled uniformly. In a sequence of eight papers between 1959-1968 they laid the foundation of Random Graph theory

What are we after?

 $ER_n(p)$ with n = 1000

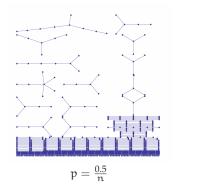


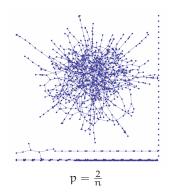


Pic source: van der Hofstad (2017)

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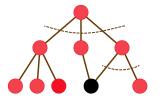
 \succ If $p = \frac{\lambda}{n}$, then there is *phase transition around* $\lambda = 1$

- → λ < 1: All components are small
- ⇒ λ > 1: There is a unique *giant component*

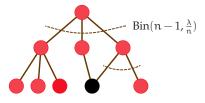
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To analyze component sizes: Gradually explore graph in BFS starting from any node, e.g., node 1

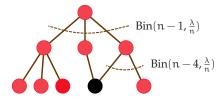
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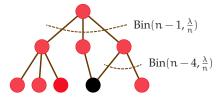
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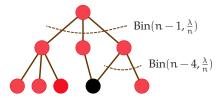


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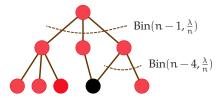
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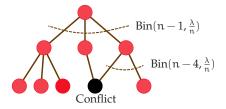


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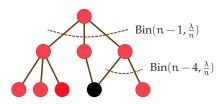
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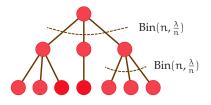
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- Depletion of vertices
- ➡ Conflicts among new vertices

Domination by Branching Process

Let's consider a random process without Depletion and Conflicts

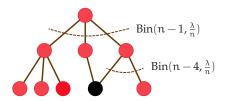


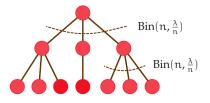


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Simple Fact

Let $N_k = \#$ vertices at depth k for $ER_n(\frac{\lambda}{n})$ exploration, and \bar{N}_k denotes same for Branching process. There is a coupling such that w.p. 1

 $N_k \leq \bar{N}_k \quad \forall k \geq 0$

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Theorem: Subcritical $ER_n(\frac{\lambda}{n})$

If $\lambda < 1$, then

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> Proof uses Large Deviation estimates for branching process survival prob

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Lets again look at upper bounding Branching Process (BP)

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 $\succ \zeta_{\lambda}$ satisfies is a positive solution of $1 - \zeta = e^{-\lambda\zeta}$

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Theorem: Supercritical $ER_n(\frac{\lambda}{n})$

Let $C_{(i)}$:= i-th largest component of $\text{ER}_n(\frac{\lambda}{n})$. If $\lambda > 1$, then as $n \to \infty$

$$\frac{C_{(1)}}{n} \xrightarrow{\mathbb{P}} \zeta_{\lambda} \quad \text{and} \quad \frac{C_{(2)}}{n} \xrightarrow{\mathbb{P}} 0$$

$$A \text{ unique giant component emerges..}$$

van der Hofstad: Random Graph and Complex Networks, Vol 2

Lemma 1: First moment

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Next, prove two lemmas but before that...

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Next, lets prove two lemmas

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Now,

 $\mathbb{P}(C(\mathfrak{u}) \geqslant L)$

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 $\mathbb{P}(\mathsf{C}(\mathsf{u}) \geqslant \mathsf{L}) \approx \mathbb{P}(\mathsf{BP} \geqslant \mathsf{L}) \qquad \qquad (\text{Exploration of } \mathsf{u} = \mathsf{BP} \text{ w.p.} \approx 1)$

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Key fact 1: Local neighborhood of u is approximately BP whp

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Key fact 1: Local neighborhood of u is approximately BP whp

> Theory of approximating local neighborhood of graphs is called Local-weak convergence (Christian's talk)

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Take two uniform vertices u_1, u_2 . The second term equals

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Enough to show:

$$\lim_{L \to \infty} \limsup_{n \to \infty} \mathbb{P} \big(C(u_1) \ge L, C(u_2) \ge L, u_1 \not\leftrightarrow u_2 \big) = 0$$

Two large components cannot be disjoint...

To show $\mathbb{P}(C(u_1) \ge L, C(u_2) \ge L, u_1 \nleftrightarrow u_2) \approx 0$, suffices to prove $\mathbb{P}(u_1 \nleftrightarrow u_2 \mid C(u_1) \ge L, C(u_2) \ge L) \approx 0$

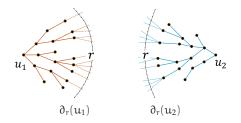
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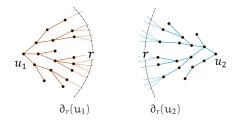
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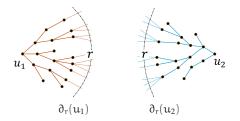


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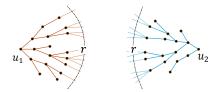
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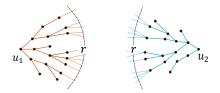


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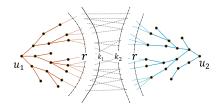
➤ Advantage with conditioning on $\{\partial_r(u_1), \partial_r(u_2) \neq \varnothing\}$ is we can now explore rest of the graph



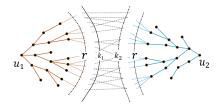
▷ Condition on $\{\partial_r(u_1), \partial_r(u_2) \neq \emptyset\}$ for large r



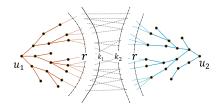
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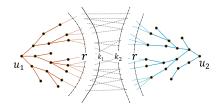


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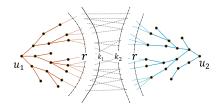
 $\mathbb{P}(\text{no edge between boundaries}) = \left(1 - \frac{\lambda}{n}\right)^{\omega_{\pi}\sqrt{n} \times \sqrt{n}} \approx e^{-\lambda \omega_{\pi}} \to 0$



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Takeaway: If there are two components with large boundary, we can grow them until boundary has size \sqrt{n} and then they intersect

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Key fact 2: $\lim_{L \to \infty} \limsup_{n \to \infty} \mathbb{P} (C(u_1) \ge L, C(u_2) \ge L, u_1 \not\leftrightarrow u_2) = 0$

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Let $C_{(i)}$:= i-th largest component of $ER_n(\frac{\lambda}{n})$. If $\lambda > 1$, then

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➤ van der Hofstad (2021) proved this for general graphs that converge in local-weak convergence sense

Before moving on to other models, lets see another useful application of the above ideas...

Typical distances in Erdős-Rényi

Typical distance: Graph distance between two uniform vertices u₁, u₂

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Let $\lambda > 1$. Conditionally on u_1 , u_2 in same component (i.e., dist $(u_1, u_2) \neq \infty$)

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0 Keep growing neighborhoods from u_1 , u_2 . Recall $\mathbb{E}[\text{#conflicts}] \leq \lambda^{2k - \log_{\lambda} n}$

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- **2** Around $k = \frac{1}{2} \log_{\lambda} n + \omega_n$, neighborhoods start intersecting

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$$\frac{\operatorname{dist}(\mathfrak{u}_1,\mathfrak{u}_2)}{\log_{\lambda} \mathfrak{n}} \xrightarrow{\mathbb{P}} 1$$

Proof: Again use neighborhood growth idea...

0 Keep growing neighborhoods from u_1 , u_2 . Recall $\mathbb{E}[\text{#conflicts}] \leq \lambda^{2k - \log_{\lambda} n}$

- → They are disjoint until boundary sizes become \sqrt{n} , i.e., $k \lesssim \frac{1}{2} \log_{\lambda} n$
- Shortest path correspond to first intersection of neighborhoods

2 Around
$$k = \frac{1}{2} \log_{\lambda} n + \omega_{n}$$
, neighborhoods start intersecting
dist $(u_{1}, u_{2}) = \frac{1}{2} \log_{\lambda} n + \frac{1}{2} \log_{\lambda} n + o(\log_{\lambda} n) = \log_{\lambda} n + o(\log_{\lambda} n)$